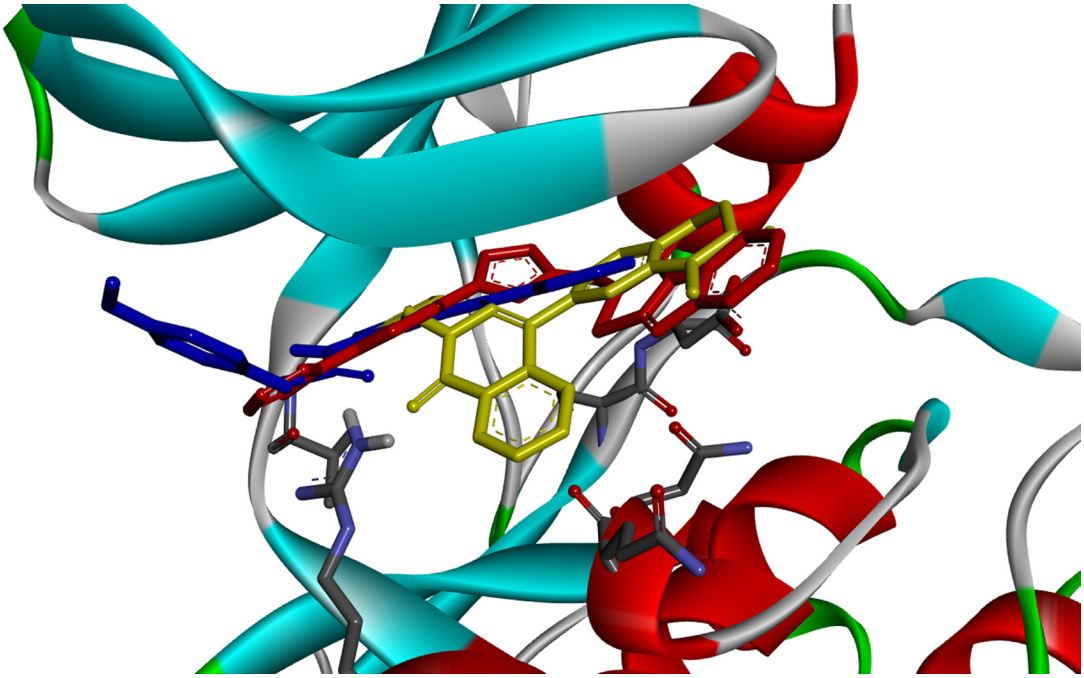
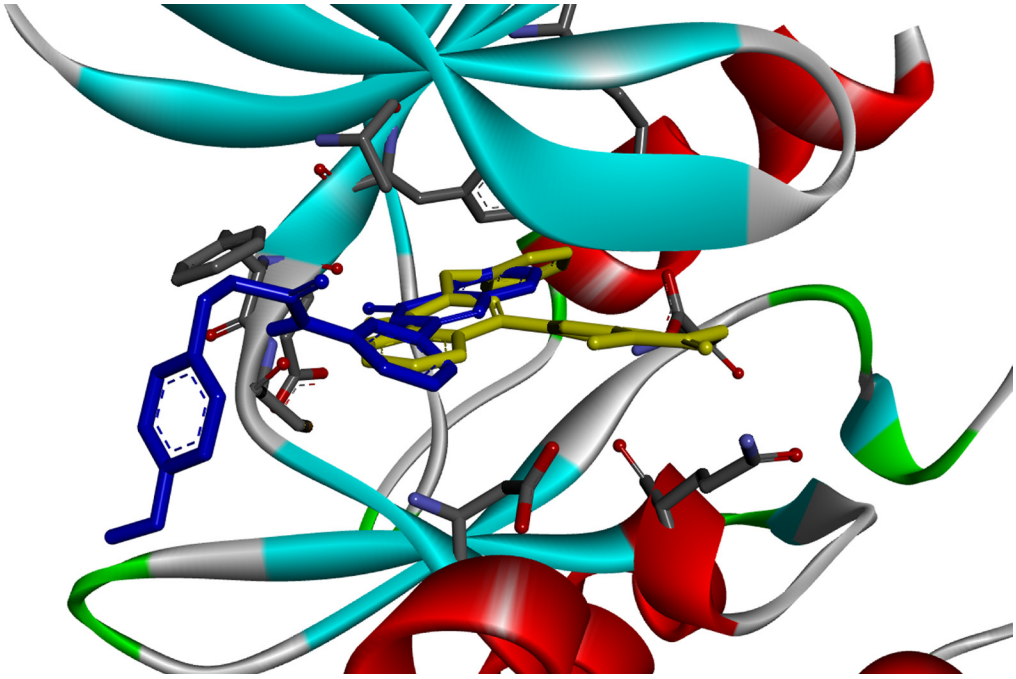
Designing of dual inhibitors for GSK-3*β* and CDK5: Virtual screening and *in vitro* biological activities study

SUPPLEMENTARY FIGURES AND TABLE



Supplementary Figure 1: The binding conformations of Compound 6, Compound 8 and Compound 9 for GSK-3β.



Supplementary Figure 2: The binding conformations of Compound 8 and Compound 9 for CDK5.

Supplementary Table 1: Docking energies of 127 compounds

See Supplementary File 1